8-579 ACCESS DB # 198152

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Scientific and Technical Information Center

SEARCH REQUEST FORM

Requester's Full Name: S. Kumar Examiner #: 69594 Date: 8/10/06

Art Unit: 1621 Phone Number: 2-0640 Serial Number: 10/528668

Location (Bldg/Room#): REM (Mailbox #): 5C18 Results Format Preferred (circle): PAPER DISK

5C03 MG

****

To ensure an efficient and quality search, please attach a copy of the cover sheet, claims, and abstract or fill out the following:

Title of Invention: Propargylether derivatives, a process for their preparation....

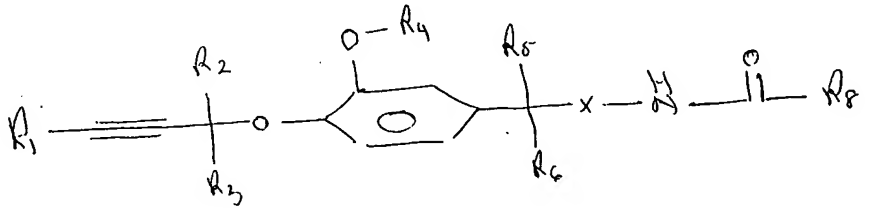
Inventors (please provide full names): Clemens Lamberth et. al.

Earliest Priority Date: 10/10/02

Search Topic:

Please provide a detailed statement of the search topic, and describe as specifically as possible the subject matter to be searched. Include the elected species or structures, keywords, synonyms, acronyms, and registry numbers, and combine with the concept or utility of the invention. Define any terms that may have a special meaning. Give examples or relevant citations, authors, etc., if known.

For Sequence Searches Only *Please include all pertinent information (parent, child, divisional, or issued patent numbers) along with the appropriate serial number.*

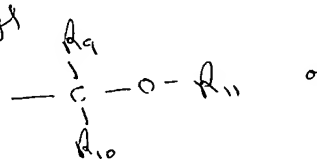
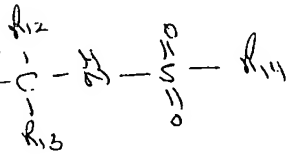


X is O or NR_7

R_1 is H, alkyl etc.

R_2-R_3, R_5-R_7 are H, alkyl etc.

R_4 is alkyl

R_8 is  or 

R_9 is Phenyl, naphthyl etc.

R_{10}, R_{11} H, alkyl etc.

R_{12} alkyl etc.

R_{13} H, alkyl etc.

R_{14} alkyl etc.

****

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Searcher: Jan

Searcher Phone #: 22504

Searcher Location:

Date Searcher Picked Up: 8/14/06

Date Completed: 8/14/06

Searcher Prep & Review Time: 15

Online Time: 110

Type of Search

____ NA Sequence (#)

____ AA Sequence (#)

✓ Structure (#)

____ Bibliographic

____ Litigation

____ Fulltext

____ Other

Vendors and cost where applicable

✓ STN ____ Dialog

____ Questel/Orbit ____ Lexis/Nexis

____ Westlaw ____ WWW/Internet

____ In-house sequence systems

____ Commercial ____ Oligomer ____ Score/Length

____ Interference ____ SPDI ____ Encode/Transl

____ Other (specify)

STIC Search Report
Biotech-Chem Library

STIC Database Tracking Number: 198152

TO:	From:
Shailendra Kumar Location: 5c03 / 5c18 Monday, August 14, 2006 Art Unit: 1621 Phone: 571-272-0640 Serial Number: 10 / 528668	Jan Delaval Location: Biotech-Chem Library Remsen 1a51 Phone: 571-272-2504 jan.delaval@uspto.gov

Search Notes

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FILE 'REGISTRY' ENTERED AT 07:53:02 ON 14 AUG 2006
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DICTIONARY FILE UPDATES:  11 AUG 2006  HIGHEST RN 900864-99-5

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http://www.cas.org/ONLINE/UG/regprops.html

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L12             STR
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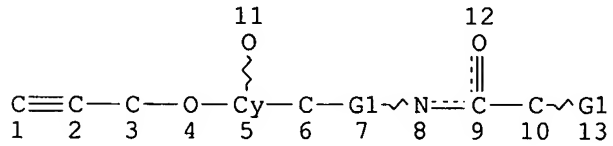


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NODE ATTRIBUTES:
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS  13

STEREO ATTRIBUTES: NONE
L14          54 SEA FILE=REGISTRY SSS FUL L12

100.0% PROCESSED     320 ITERATIONS                    54 ANSWERS
SEARCH TIME: 00.00.01


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               E LAMBERTH/AU
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L2            68 S E5,E8
                 E ZELLER/AU
L3             5 S E3
                 E ZELLER M/AU
L4           167 S E3-E7,E17,E18
                 E SYNGENTA/PA,CS
L5          1218 S E3,E4 OR SYNGENTA?/PA,CS
                 SEL RN L1

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L9               STR L7
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L19            1 S L18 AND L1-L5

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L20            0 S L14

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FILE 'HCAPLUS' ENTERED AT 07:53:11 ON 14 AUG 2006
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FILE LAST UPDATED: 13 Aug 2006  (20060813/ED)

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 This file contains CAS Registry Numbers for easy and accurate
 substance identification.
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=> d l19 bib abs hitstr retable

L19 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2006 ACS on STN
AN 2004:333686 HCAPLUS
DN 140:357056
TI Preparation of novel propargyl ether derivatives for controlling phytopathogenic microorganisms
IN **Lamberth, Clemens; Zeller, Martin**
PA **Syngenta Participations Ag, Switz.**
SO PCT Int. Appl., 57 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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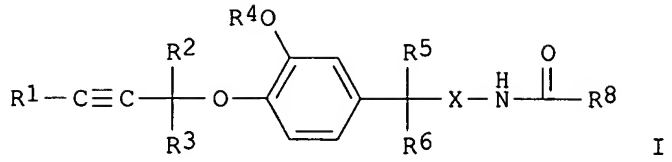

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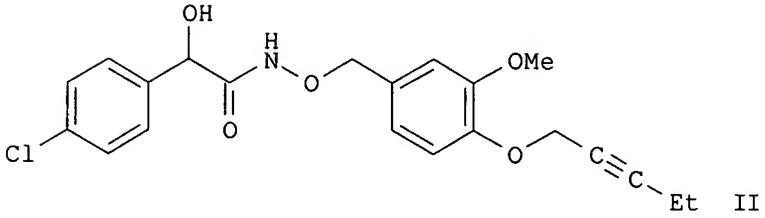

II

AB The title compds. [I; R1= H, (un)substituted alkyl, cycloalkyl, aryl; R2, R3, R5-R7 = H, alkyl; R4 = (un)substituted alkyl; X = O, NR7; R8 = CR9R10OR11, CR12R13NHSO2R14 (wherein R9 = (un)substituted (hetero)aryl; R10, R11 = H, (un)substituted alkyl, alkenyl, alkynyl; R12 = (un)substituted alkyl, cycloalkyl, aryl, heteroaryl; R13 = H, (un)substituted alkyl, alkenyl or alkynyl; R14 = (un)substituted alkyl, NH2)] which possess plant protecting properties and may advantageously be employed in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi, were prepared E.g, a multi-step synthesis of II, starting from 4-hydroxymethyl-2-methoxyphenol and MeCH2C.tplbond.CCH2OH, was given. Representative compds. I showed at least 80% inhibition of fungal infestation in 3 biol. tests.

IT **681434-45-7P 681434-46-8P 681434-47-9P**
681434-48-0P 681434-49-1P 681434-50-4P
681434-51-5P 681434-52-6P 681434-53-7P
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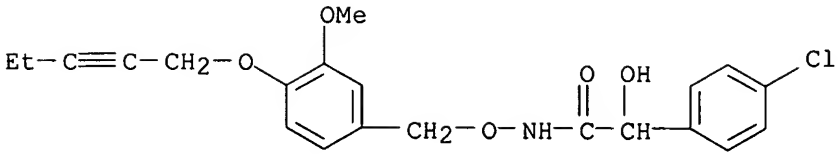
681435-03-0P 681435-04-1P 681435-05-2P
RL: AGR (Agricultural use); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of novel propargyl ether derivs. for controlling phytopathogenic microorganisms)

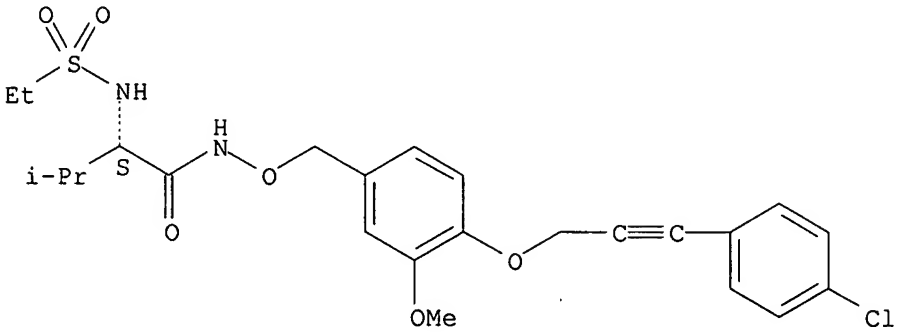
RN 681434-45-7 HCAPLUS

CN Benzeneacetamide, 4-chloro-α-hydroxy-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



RN 681434-46-8 HCAPLUS
CN Butanamide, N-[[4-[[3-(4-chlorophenyl)-2-propynyl]oxy]-3-methoxyphenyl]methoxy]-2-[(ethylsulfonyl)amino]-3-methyl-, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 681434-47-9 HCAPLUS
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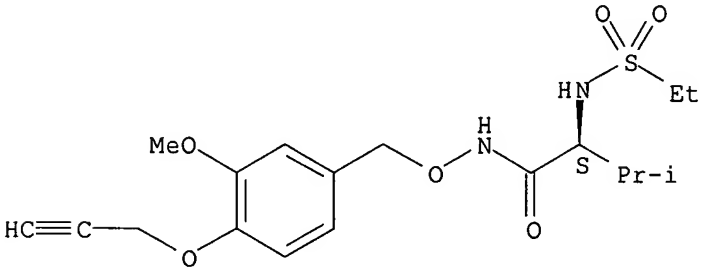
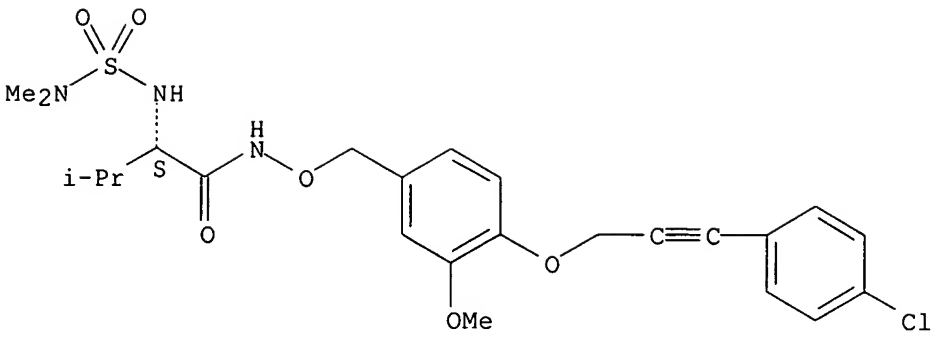
Absolute stereochemistry.



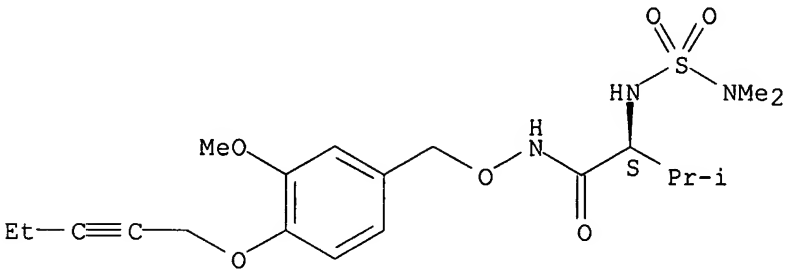
RN 681434-48-0 HCAPLUS
CN Butanamide, N-[[4-[[3-(4-chlorophenyl)-2-propynyl]oxy]-3-methoxyphenyl]methoxy]-2-[[(dimethylamino)sulfonyl]amino]-3-methyl-, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 681434-49-1 HCAPLUS
CN Butanamide, 2-[[(dimethylamino)sulfonyl]amino]-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]-3-methyl-, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 681434-50-4 HCAPLUS
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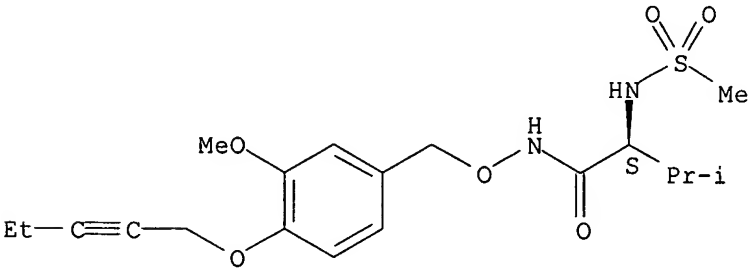
Absolute stereochemistry.



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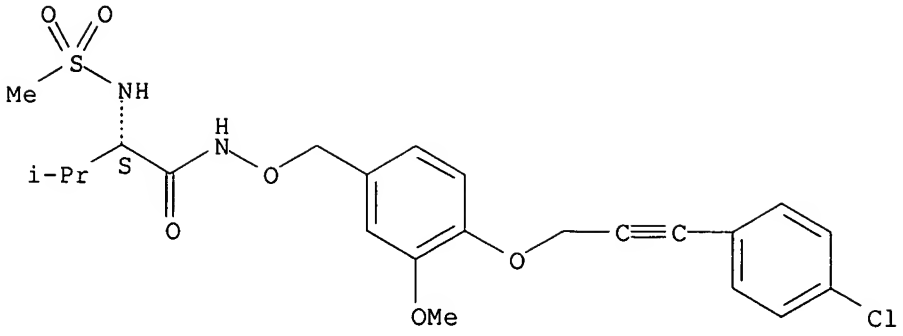
Absolute stereochemistry.



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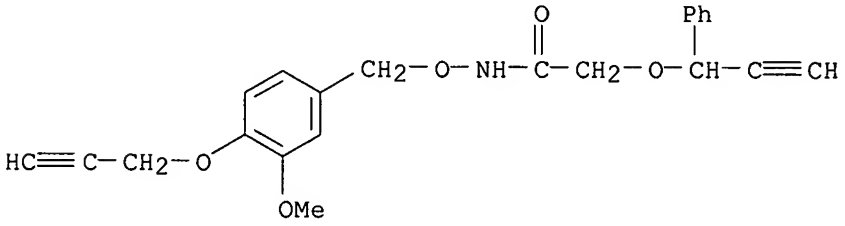

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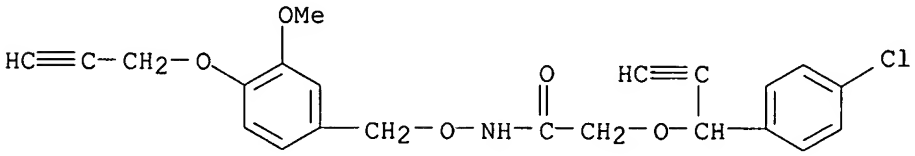

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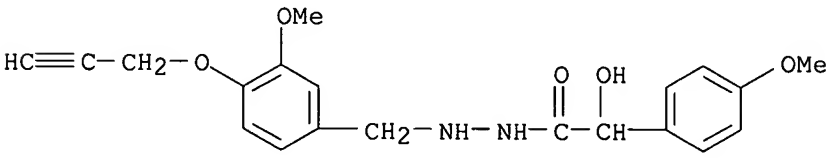

RN 681434-55-9 HCAPLUS
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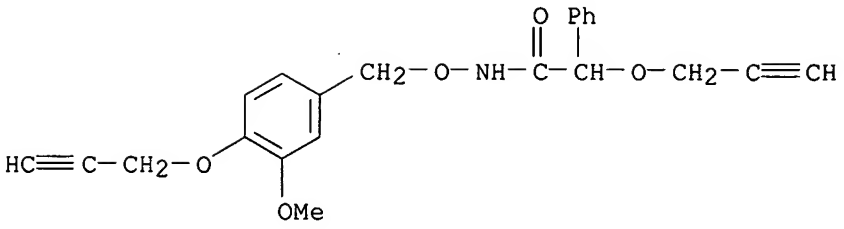

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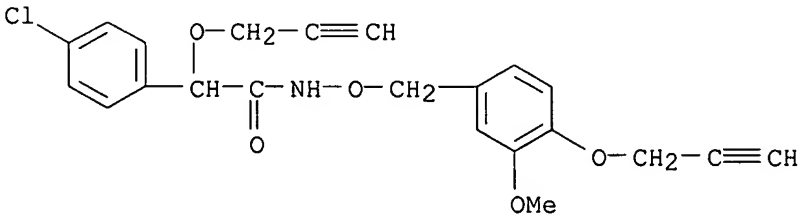

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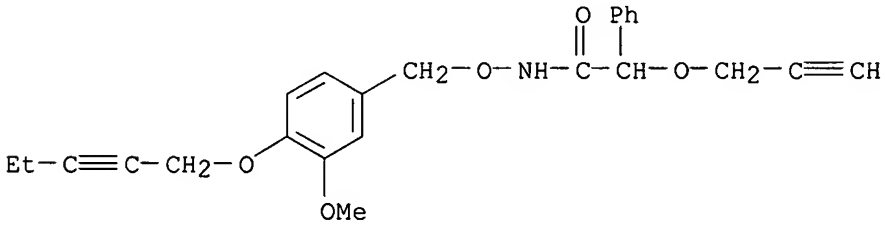

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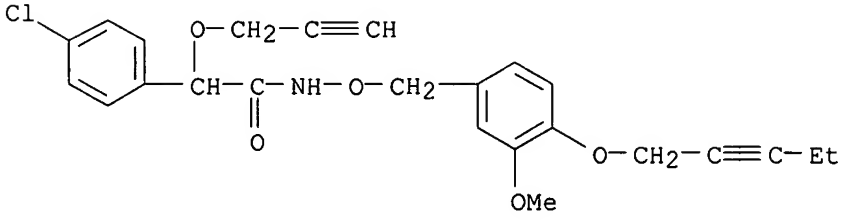

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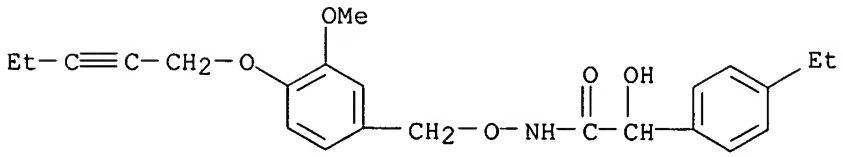

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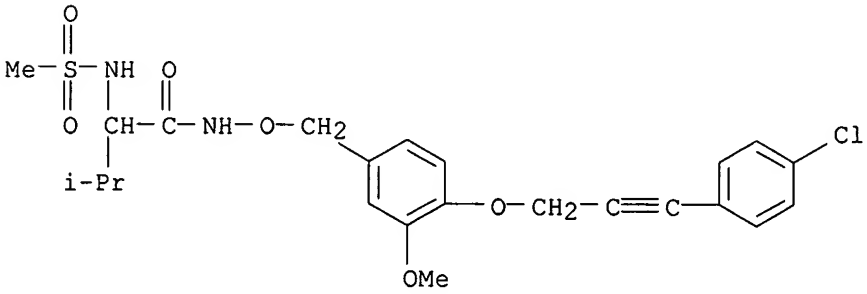

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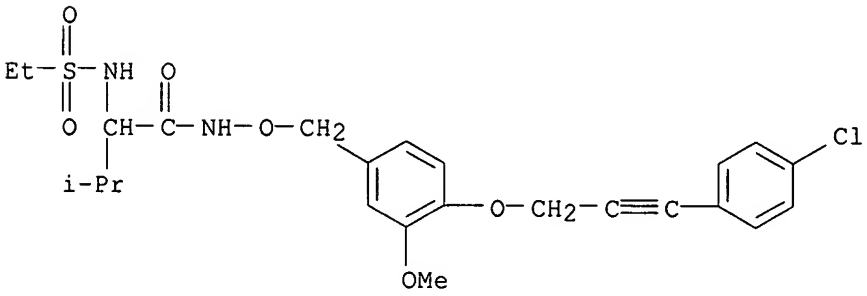
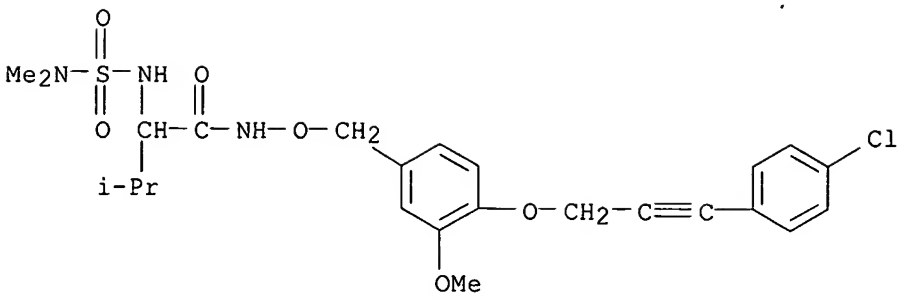
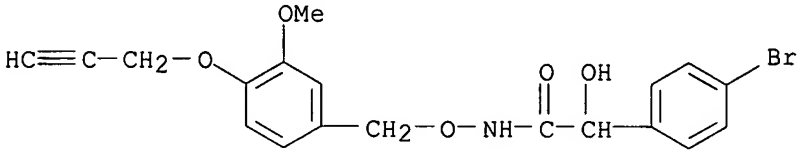


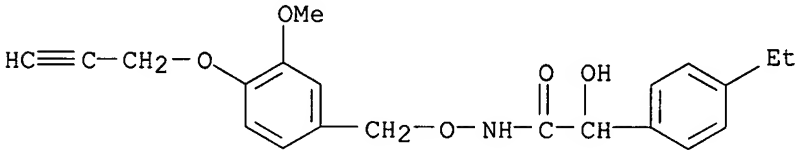
RN 681434-62-8 HCAPLUS
CN Butanamide, N-[[4-[[3-(4-chlorophenyl)-2-propynyl]oxy]-3-methoxyphenyl]methoxy]-2-[[(dimethylamino)sulfonyl]amino]-3-methyl- (9CI) (CA INDEX NAME)



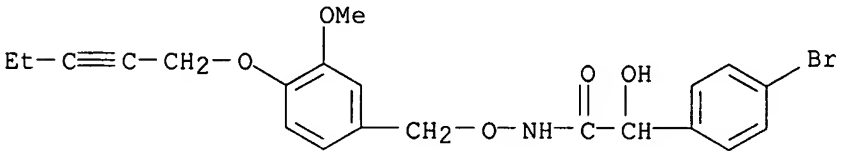
RN 681434-63-9 HCAPLUS
CN Benzeneacetamide, 4-bromo-α-hydroxy-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



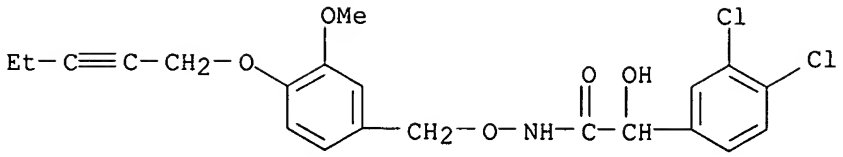
RN 681434-64-0 HCAPLUS
CN Benzeneacetamide, 4-ethyl-α-hydroxy-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



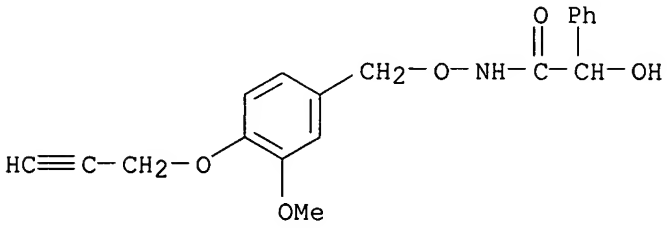
RN 681434-65-1 HCAPLUS
CN Benzeneacetamide, 4-bromo-α-hydroxy-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



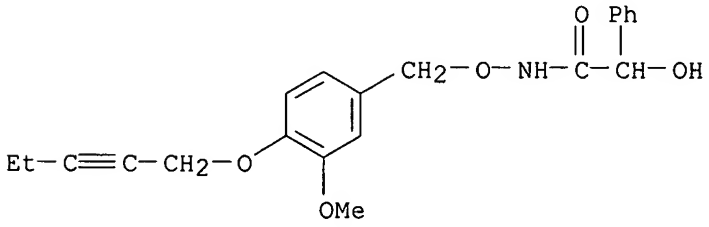
RN 681434-66-2 HCAPLUS
CN Benzeneacetamide, 3,4-dichloro-α-hydroxy-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



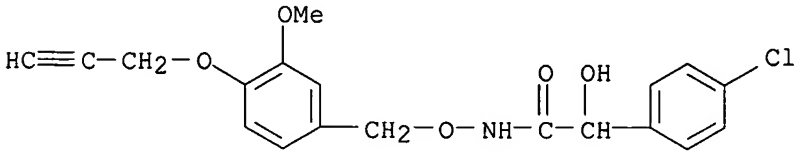
RN 681434-74-2 HCAPLUS
CN Benzeneacetamide, α-hydroxy-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



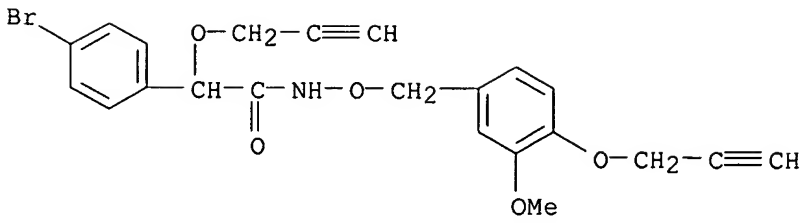
RN 681434-75-3 HCAPLUS
CN Benzeneacetamide, α-hydroxy-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



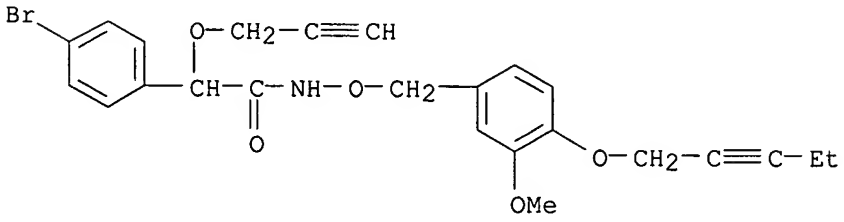
RN 681434-76-4 HCAPLUS
CN Benzeneacetamide, 4-chloro-α-hydroxy-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



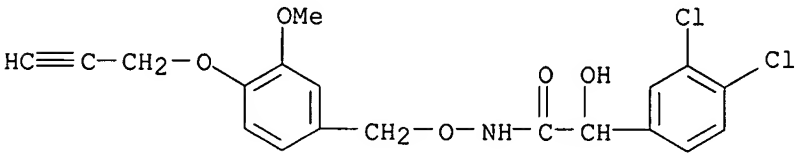
RN 681434-77-5 HCAPLUS
CN Benzeneacetamide, 4-bromo-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



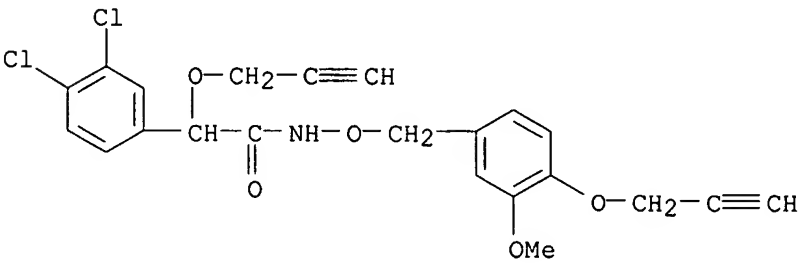
RN 681434-78-6 HCAPLUS
CN Benzeneacetamide, 4-bromo-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



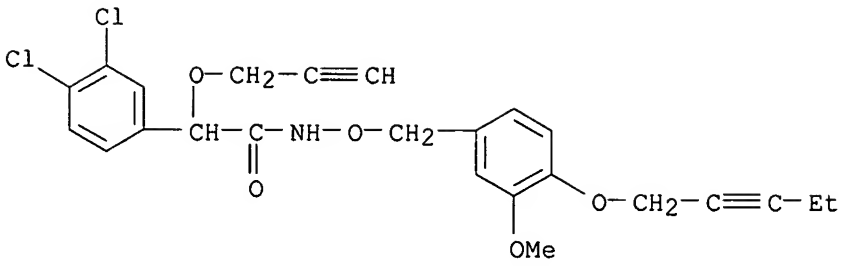
RN 681434-79-7 HCAPLUS
CN Benzeneacetamide, 3,4-dichloro-α-hydroxy-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]- (9CI) (CA INDEX NAME)



RN 681434-80-0 HCAPLUS
CN Benzeneacetamide, 3,4-dichloro-N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)

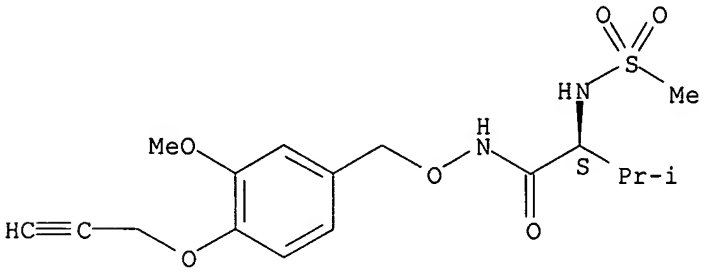


RN 681434-81-1 HCAPLUS
CN Benzeneacetamide, 3,4-dichloro-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]-α-(2-propynyloxy)- (9CI) (CA INDEX NAME)



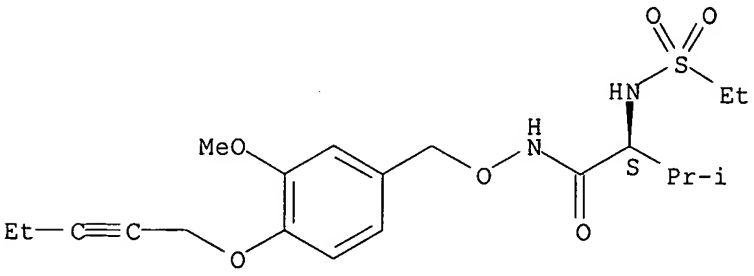
RN 681434-82-2 HCAPLUS
CN Butanamide, N-[[3-methoxy-4-(2-propynyloxy)phenyl]methoxy]-3-methyl-2-[(methylsulfonyl)amino]-, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

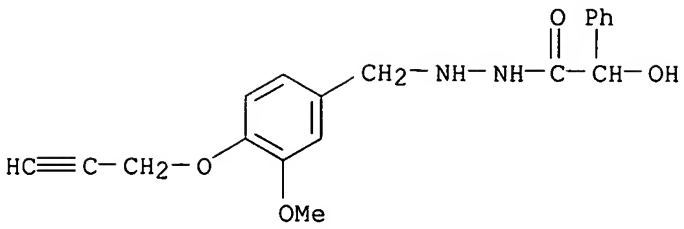


RN 681434-83-3 HCAPLUS
CN Butanamide, 2-[(ethylsulfonyl)amino]-N-[[3-methoxy-4-(2-pentynyloxy)phenyl]methoxy]-3-methyl-, (2S)- (9CI) (CA INDEX NAME)

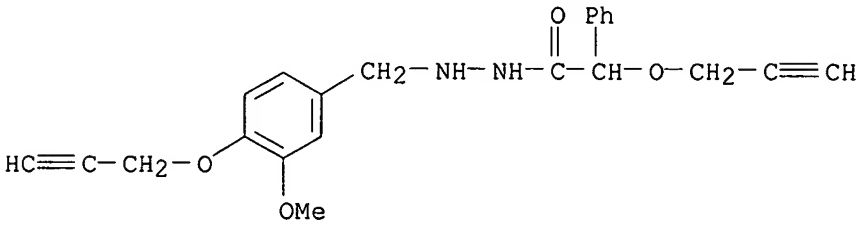
Absolute stereochemistry.



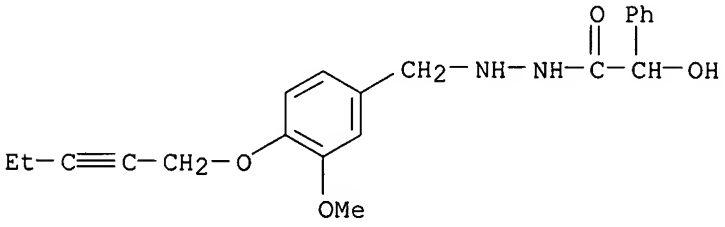
RN 681434-84-4 HCAPLUS
CN Benzeneacetic acid, α-hydroxy-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



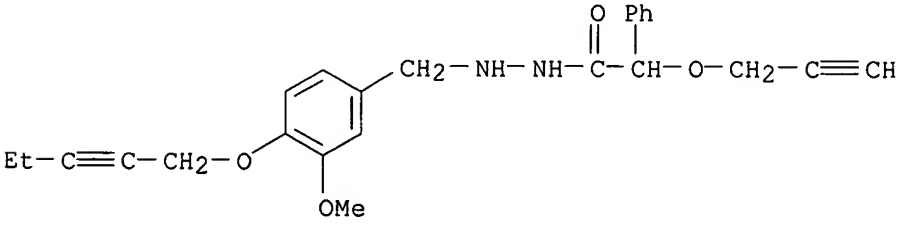
RN 681434-85-5 HCAPLUS
CN Benzeneacetic acid, α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



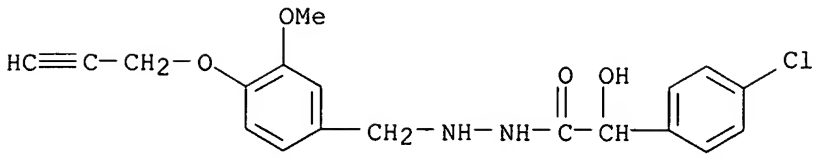
RN 681434-86-6 HCAPLUS
CN Benzeneacetic acid, α-hydroxy-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



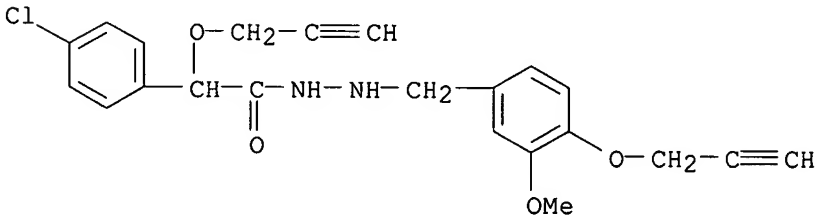
RN 681434-87-7 HCAPLUS
CN Benzeneacetic acid, α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



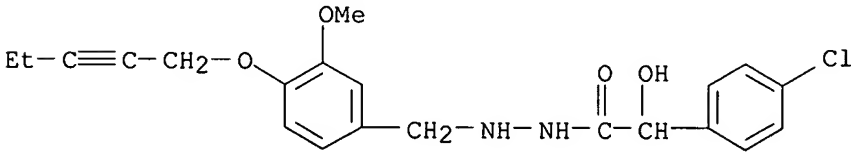
RN 681434-88-8 HCAPLUS
CN Benzeneacetic acid, 4-chloro-α-hydroxy-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



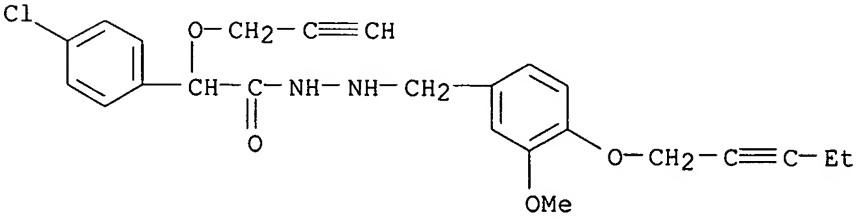
RN 681434-89-9 HCAPLUS
CN Benzeneacetic acid, 4-chloro-α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



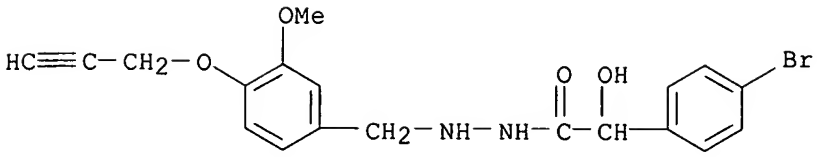
RN 681434-90-2 HCAPLUS
CN Benzeneacetic acid, 4-chloro-α-hydroxy-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



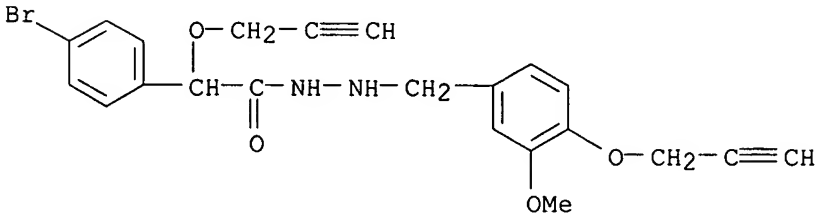
RN 681434-91-3 HCAPLUS
CN Benzeneacetic acid, 4-chloro-α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



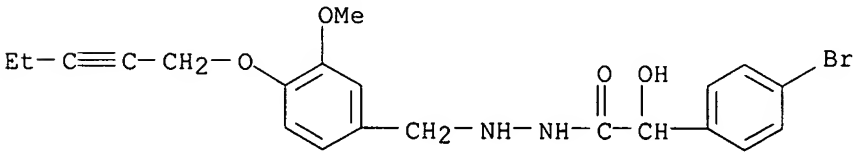
RN 681434-92-4 HCAPLUS
CN Benzeneacetic acid, 4-bromo-α-hydroxy-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



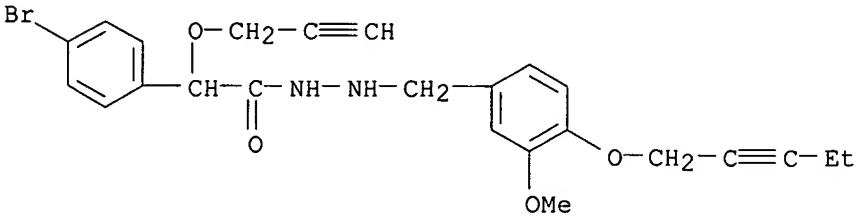
RN 681434-93-5 HCAPLUS
CN Benzeneacetic acid, 4-bromo-α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



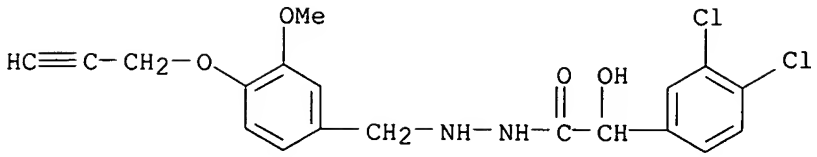
RN 681434-94-6 HCAPLUS
CN Benzeneacetic acid, 4-bromo-α-hydroxy-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



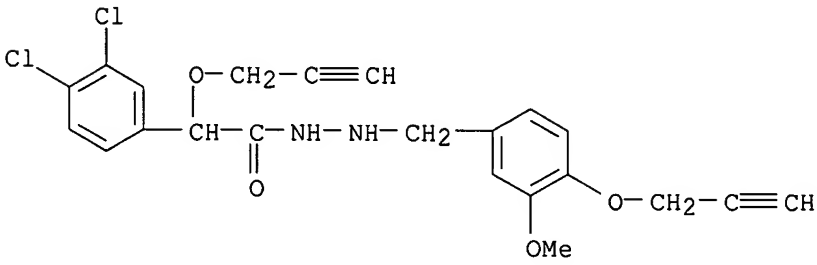
RN 681434-95-7 HCAPLUS
CN Benzeneacetic acid, 4-bromo-α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



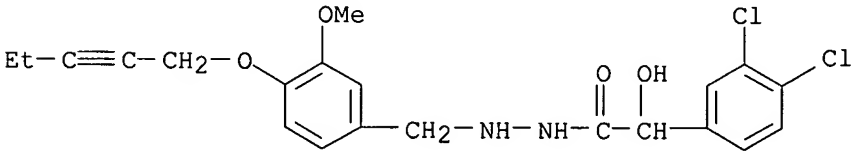
RN 681434-96-8 HCAPLUS
CN Benzeneacetic acid, 3,4-dichloro-α-hydroxy-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



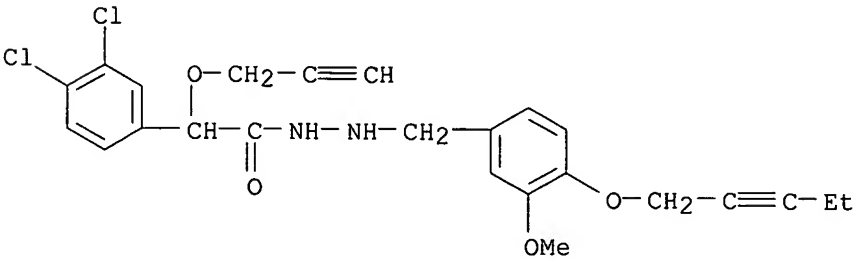
RN 681434-97-9 HCAPLUS
CN Benzeneacetic acid, 3,4-dichloro-α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



RN 681434-98-0 HCAPLUS
CN Benzeneacetic acid, 3,4-dichloro-α-hydroxy-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

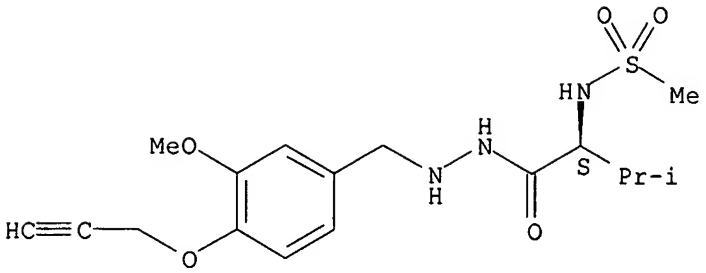


RN 681434-99-1 HCAPLUS
CN Benzeneacetic acid, 3,4-dichloro-α-(2-propynyloxy)-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)



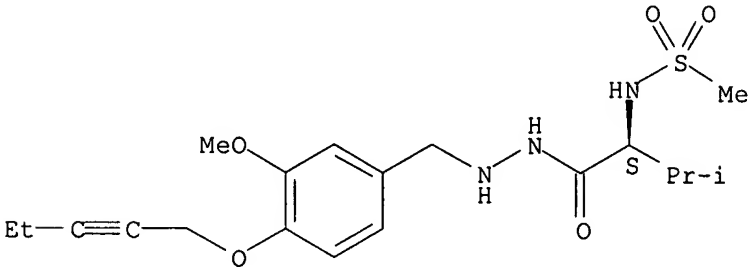
RN 681435-00-7 HCAPLUS
CN L-Valine, N-(methylsulfonyl)-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

Absolute stereochemistry.



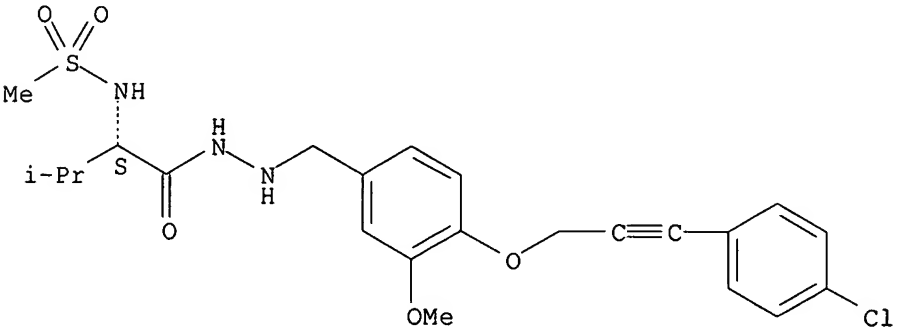
RN 681435-01-8 HCAPLUS
CN L-Valine, N-(methylsulfonyl)-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

Absolute stereochemistry.



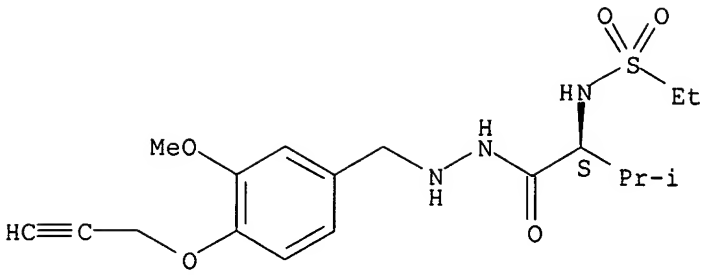
RN 681435-02-9 HCAPLUS
CN L-Valine, N-(methylsulfonyl)-, 2-[[4-[[3-(4-chlorophenyl)-2-propynyl]oxy]-3-methoxyphenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

Absolute stereochemistry.



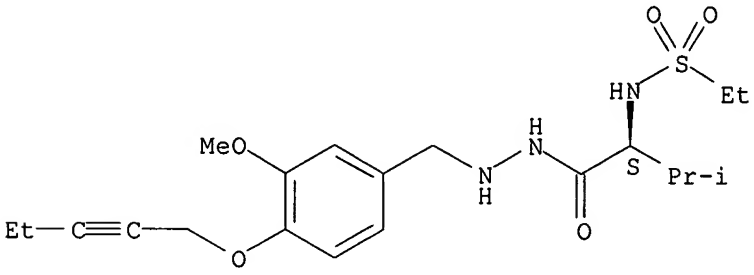
RN 681435-03-0 HCAPLUS
CN L-Valine, N-(ethylsulfonyl)-, 2-[[3-methoxy-4-(2-propynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

Absolute stereochemistry.



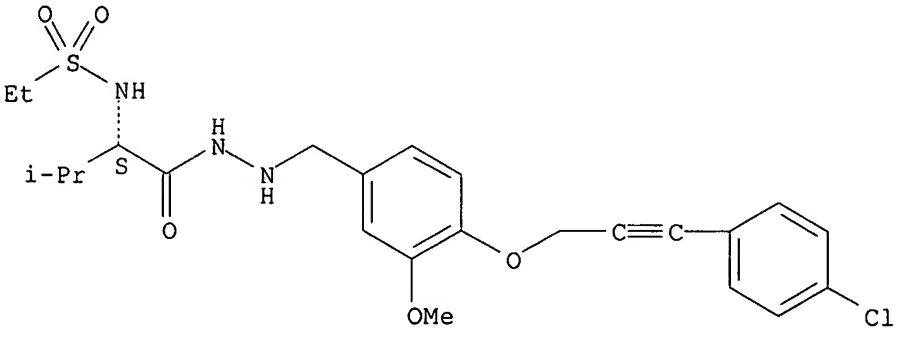
RN 681435-04-1 HCAPLUS
CN L-Valine, N-(ethylsulfonyl)-, 2-[[3-methoxy-4-(2-pentynyloxy)phenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 681435-05-2 HCAPLUS
CN L-Valine, N-(ethylsulfonyl)-, 2-[[4-[[3-(4-chlorophenyl)-2-propynyl]oxy]-3-methoxyphenyl]methyl]hydrazide (9CI) (CA INDEX NAME)

Absolute stereochemistry.



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